

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
 Data File : BF096438.D
 Acq On : 3 Jul 2017 13:51
 Operator : SJ/MA
 Sample : I3908-08 5X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-59-1.5-3

Manual Integrations
 APPROVED

mohammad
 7/5/2017 11:41:10 AM

Quant Time: Jul 03 16:49:52 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 01:43:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	117455	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	423894	20.00	ng	-0.02
38) Acenaphthene-d10	9.76	164	171436	20.00	ng	-0.01
63) Phenanthrene-d10	11.23	188	237861	20.00	ng	-0.02
75) Chrysene-d12	13.86	240	187849	20.00	ng	-0.02
86) Perylene-d12	15.27	264	160362	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	228328	28.69	ng	-0.02
7) Phenol-d6	6.35	99	293764	30.03	ng	-0.02
23) Nitrobenzene-d5	7.27	82	139015	19.71	ng	-0.03
41) 2,4,6-Tribromophenol	10.54	330	36686	27.40	ng	-0.02
44) 2-Fluorobiphenyl	9.08	172	260234	25.95	ng	-0.02
78) Terphenyl-d14	12.82	244	137200	15.61	ng	-0.02
Target Compounds						Qvalue
31) Naphthalene	8.03	128	47376	2.367	ng	99
37) 2-Methylnaphthalene	8.72	142	35836	2.813	ng	95
49) Dimethylphthalate	9.47	163	30517	2.385	ng	97
74) Fluoranthene	12.44	202	106657	8.460	ng	98
77) Pyrene	12.67	202	103123	6.839	ng	97
80) Benzo(a)anthracene	13.86	228	61474	5.651	ng	# 91
82) Chrysene	13.89	228	63017m	5.532	ng	
85) Indeno(1,2,3-cd)pyrene	16.63	276	23713	2.637	ng	97
87) Benzo(b)fluoranthene	14.87	252	70465m	7.013	ng	
88) Benzo(k)fluoranthene	14.90	252	25008m	2.782	ng	
89) Benzo(a)pyrene	15.22	252	42962	4.789	ng	# 90
91) Benzo(g,h,i)perylene	17.05	276	22757	3.112	ng	# 87

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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